10/551,771 08/28/2006

=> d his ful

(FILE 'HOME' ENTERED AT 11:10:56 ON 28 AUG 2006)

FILE 'REGISTRY' ENTERED AT 11:11:34 ON 28 AUG 2006

L1 STRUCTURE UPLOADED L2

STRUCTURE UPLOADED

D.L1

L*** DEL 1676 DL 2 D L2

50 SEA SSS SAM L1 L3 50 SEA SSS SAM L2 L4

T.5 57174 SEA SSS FUL L1 L6 413178 SEA SSS FUL L2

L7 228236 SEA SUB=L6 SSS FUL L1

FILE 'HCAPLUS' ENTERED AT 11:14:08 ON 28 AUG 2006 L8 22139 SEA PLU=ON L7

FILE 'REGISTRY' ENTERED AT 11:15:23 ON 28 AUG 2006

STRUCTURE UPLOADED L9

D L9

13 SEA SSS SAM L9 L10

L11 18 SEA SSS FUL L9 AND L2

FILE 'HCAPLUS' ENTERED AT 11:16:22 ON 28 AUG 2006

L12 13 SEA PLU=ON L11

D L12 1-13 IBIB HITSTR

FILE HOME

FILE REGISTRY

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 27 AUG 2006 HIGHEST RN 904741-41-9 DICTIONARY FILE UPDATES: 27 AUG 2006 HIGHEST RN 904741-41-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

FILE HCAPLUS

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10/551,771 08/28/2006

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FILE COVERS 1907 - 28 Aug 2006 VOL 145 ISS 10 FILE LAST UPDATED: 27 Aug 2006 (20060827/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> d que sta

L2 STR

Structure attributes must be viewed using STN Express query preparation. L9 $$\operatorname{\mathtt{STR}}$$

$$CH_2$$
 OH CH_2 OH

Structure attributes must be viewed using STN Express query preparation.

L11 18 SEA FILE=REGISTRY SSS FUL L9 AND L2

L12 13 SEA FILE=HCAPLUS PLU=ON L11

C:\Program Files\Stnexp\Queries\hugh1.str

chain nodes :
 1 2 3 4 5 6 7 10 11 12 13 14 17 18 19
chain bonds :
 1-2 2-3 3-4 4-5 4-7 5-6 6-10 10-11 11-12 11-14 12-13 13-17
 17-18 18-19
exact/norm bonds :
 1-2 2-3 3-4 4-5 4-7 5-6 6-10 10-11 11-12 11-14 12-13 13-17
 17-18 18-19

Match level:
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 17:CLASS 18:CLASS 19:CLASS

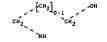
C:\Program Files\Stnexp\Queries\hugh2.str

chain nodes : 1 2 3 4 7 chain bonds : 1-2 2-3 3-4 4-7

exact/norm bonds : 1-2 2-3 3-4 4-7

Match level: 1:CLASS 2:CLASS 3:CLASS 4:CLASS 7:CLASS

C:\Program Files\Stnexp\Queries\hugh3.str



2-(33)0-1

chain nodes :
 1 2 3 4 7
chain bonds :
 1-2 2-3 3-4 4-7
exact/norm bonds :

1-2 2-3 3-4 4-7

Match level: 1:CLASS 2:CLASS 3:CLASS 4:CLASS 7:CLASS

> d his ful

(FILE 'HOME' ENTERED AT 11:27:55 ON 28 AUG 2006)

FILE 'REGISTRY' ENTERED AT 11:27:59 ON 28 AUG 2006 Ll STRUCTURE UPLOADED D STRUCTURE UPLOADED L2D L3 0 SEA SSS SAM L2 L40 SEA SSS FUL L2 L5 STRUCTURE UPLOADED L6 0 SEA SSS FUL L5 L7 STRUCTURE UPLOADED L8 0 SEA SSS FUL L7

FILE HOME

FILE REGISTRY

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 27 AUG 2006 HIGHEST RN 904741-41-9 DICTIONARY FILE UPDATES: 27 AUG 2006 HIGHEST RN 904741-41-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=> d que sta

L7 STR

Structure attributes must be viewed using STN Express query preparation. L8 $\,$ 0 SEA FILE=REGISTRY SSS FUL L7 $\,$

100.0% PROCESSED 700042 ITERATIONS SEARCH TIME: 00.00.41

0.ANSWERS

```
chain nodes :
   1 2 3 4 5 6 7 8 9 10 11 12 13
                                        14
                                            15
                                               16 17
                                                     18
                                                                 21
   22 23 24 25 26 27 28 29
                              30 31 32 33 34 35 36
   1-2 1-29 1-36 2-3 2-4 3-5 5-6 5-24
                                       6-7 6-14 8-10 8-9 8-23
   10-11 11-12 11-28 12-13 14-15 15-16 15-25 16-17 16-18 18-19
                                                                19-20
   19-26 20-21 20-22 21-23 23-27 24-31 25-32 26-33 27-34 28-35
exact/norm bonds :
                          3-5 5-6 5-24
   1-2 1-29 1-36
                 2-3
                     2 - 4
                                        6-7
                                            6-14 8-10 8-9 8-23
                     12-13 14-15 15-16
                                       15-25
                                             16-17 16-18 18-19
   10-11
         11-12
              11-28
                                                                19-20
                     21-23 23-27 24-31
   19-26 20-21 20-22
                                       25-32
                                             26-33
                                                    27-34
                                                          28-35
                                                                29-30
```

Match level:
 1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS 36:CLASS Generic attributes:

24:

Type of chain : Linear
Saturation : Saturated
Number of Carbon Atoms : less than 7

25:

Type of chain : Linear Saturation : Saturated Number of Carbon Atoms : less than 7

26:

Type of chain : Linear Saturation : Saturated Number of Carbon Atoms : less than 7

27:

Type of chain : Linear

Saturation : Saturated Number of Carbon Atoms : less than 7

28:

Type of chain : Linear
Saturation : Saturated
Number of Carbon Atoms : less than 7

29:

Type of chain : Linear Saturation : Saturated Number of Carbon Atoms : less than 7

25:

26:

27:

Type of chain

Type of chain

Type of chain

Saturation

Saturation

```
chain nodes :
   1 2 3 4 5 6 7 8 9 10 11 12 13 14
                                                15 16
                                                        17 18 19 20
                                                                       21
   22 23 24 25 26 27 28
                             29
                                 30 31 32 33 34 35
                                                       36
   1-2 1-29 1-36 2-3 2-4 3-5 5-6 5-24
                                           6-7
                                                 6-14 8-10 8-9 8-23
   10-11 11-12 11-28 12-13 14-15 15-16 15-25 16-17 16-18 18-19
   19-26 20-21 20-22 21-23 23-27 24-31
                                           25-32 26-33 27-34
                                                               28-35
exact/norm bonds :
                       2-4 3-5 5-6 5-24 6-7
12-13 14-15 15-16 15-25
                                           6-7 6-14 8-10 8-9 8-23
15-25 16-17 16-18 18-19
                   2-3
   1-2 1-29 1-36
          11-12 11-28
                                                                      19-20
               20-22
   19-26
          20-21
                       21-23 23-27 24-31
                                           25-32
                                                  26-33 27-34
                                                               28-35
Match level :
   1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS
                                                               8:CLASS
   9:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS
                     18:CLASS 19:CLASS 20:CLASS 21:CLASS 22:CLASS
   16:CLASS
            17:CLASS
   23:CLASS
             24:CLASS
                     25:CLASS 26:CLASS
                                                  28:CLASS 29:CLASS
                                         27:CLASS
   30:CLASS
            31:CLASS
                     32:CLASS 33:CLASS
                                         34:CLASS
                                                  35:CLASS 36:CLASS
Generic attributes :
   24:
   Type of chain
                         : Linear
                         : Saturated
   Saturation
   Number of Carbon Atoms : less than 7
```

: Linear

: Linear

: Linear

Number of Carbon Atoms : less than 7

Number of Carbon Atoms : less than 7

: Saturated

: Saturated

'Saturation : Saturated Number of Carbon Atoms : less than 7

28:

Type of chain : Linear
Saturation : Saturated
Number of Carbon Atoms : less than 7

29:

Type of chain : Linear Saturation : Saturated Number of Carbon Atoms : less than 7

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=> d 112 1-13 ibib hitstr

L12 ANSWER 1 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:1290320 HCAPLUS

DOCUMENT NUMBER: 144:192568

TITLE: Amphiphilic Poly(L-lactide)-b-dendritic

Poly(L-lysine)s Synthesized with A Metal-Free Catalyst and New Dendron Initiators: Chemical Preparation and

Characterization

AUTHOR(S): Li, Yang; Li, Qiaobo; Li, Faxue; Zhang, Haiyun; Jia,

Lin; Yu, Jianyong; Fang, Qiang; Cao, Amin

CORPORATE SOURCE: Laboratory for Polymer Materials, Shanghai Institute

of Organic Chemistry, Chinese Academy of Sciences,

Shanghai, 200032, Peop. Rep. China

SOURCE: Biomacromolecules (2006), 7(1), 224-231

CODEN: BOMAF6; ISSN: 1525-7797

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal LANGUAGE: English

IT 875275-84-6P

RL: SPN (Synthetic preparation); PREP (Preparation)

(preparation and characterization of amphiphilic poly(L-lactide)-b-dendritic

poly(L-lysine)s synthesized with metal-free catalyst and new dendron

initiators)

RN 875275-84-6 HCAPLUS

CN 1,4-Dioxane-2,5-dione, 3,6-dimethyl-, (3S,6S)-, homopolymer,

 $2-[[N2,N6-bis(N2,N6-di-L-lysyl-L-lysyl)-L-lysyl] amino] ethyl \ ester \ (9CI) \\$

(CA INDEX NAME)

CM 1

CRN 875275-83-5

CMF C44 H91 N15 O8

PAGE 1-B

CM 2

CRN 33135-50-1 CMF (C6 H8 O4)x

CCI PMS

CM 3

CRN 4511-42-6 CMF C6 H8 O4

Absolute stereochemistry.

REFERENCE COUNT: 30 THERE ARE 30 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 2 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:1211350 HCAPLUS

DOCUMENT NUMBER: 144:108650

TITLE: Sequence-Defined Polypeptide-Polymer Conjugates

Utilizing Reversible Addition Fragmentation Transfer

Radical Polymerization

AUTHOR(S): ten Cate, Mattijs G. J.; Rettig, Hartmut; Bernhardt,

Kaj; Boerner, Hans G.

CORPORATE SOURCE: Max Planck Institute of Colloids and Interfaces, MPI

KGF Golm, Potsdam, 14424, Germany

SOURCE: Macromolecules (2005), 38(26), 10643-10649

CODEN: MAMOBX; ISSN: 0024-9297

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal LANGUAGE: English

IT 873078-15-0D, functionalized polystyrene-supported RL: RCT (Reactant); RACT (Reactant or reagent)

(sequence-defined polypeptide-polymer conjugates utilizing RAFT polymerization)

RN 873078-15-0 HCAPLUS

CN L- α -Asparagine, glycyl-L- α -aspartylglycyl-L-phenylalanyl-N-(2-hydroxyethyl)-, bis(1,1-dimethylethyl) ester (9CI) (CA INDEX NAME)

10/551,771 08/28/2006

Absolute stereochemistry.

REFERENCE COUNT: 36 THERE ARE 36 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 3 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:450920 HCAPLUS

DOCUMENT NUMBER: 142:482324

TITLE: Preparation of phenylalanine derivatives as

 δ -opioid receptor ligands

INVENTOR(S): Dolle, Roland E.

PATENT ASSIGNEE(S): USA

SOURCE: U.S. Pat. Appl. Publ., 34 pp., Cont.-in-part of U.S.

Ser. No. 719,627.

CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2005113295	Al	20050526	US 2004-991785	20041118
US 2005113294	A1	20050526	US 2003-719627	20031121
PRIORITY APPLN. INFO.:			US 2003-719627 A2	20031121
OTHER SOURCE(S):	MARPAT	142:482324	•	

IT 851883-43-7P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES

(preparation of phenylalanyl peptides as δ -opioid receptor ligands)

RN 851883-43-7 HCAPLUS

CN L-Phenylalaninamide, 4-(aminocarbonyl)-L-phenylalanyl-D-alanylglycyl-N-(2-hydroxyethyl)-N α -methyl- (9CI) (CA INDEX NAME)

10/551,771 08/28/2006

PAGE 1-B

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(Uses)

L12 ANSWER 4 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

2005:450919 HCAPLUS ACCESSION NUMBER:

DOCUMENT NUMBER: 142:482323

Preparation of phenylalanine derivatives as TITLE:

 δ -opioid receptor ligands

INVENTOR(S): Dolle, Roland E.

PATENT ASSIGNEE(S): Adolor Corporation, USA

SOURCE: U.S. Pat. Appl. Publ., 32 pp.

CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

	NO.				DATE			APPL:	ICAT	ION I	NO.		D	ATE	
	113294				2005	0526		US 2	003-	7196	27		2	0031	121
	113295				2005	0526		US 2	004-	9917	85		21	0041	118
WO 2005	051367		A1		2005	0609	1	WO 2	004-	US38	656		2	0041	118
W:	AE, AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,
	CN, CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,
	GE, GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,
	LK, LR,														
	NO, NZ,				-	-						•	•	-	
	TJ, TM,														
RW:	BW, GH,														
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OTHER SOURCE												_			
IT 851883-	• •														
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	eutic us		_											•	

10/551,771 08/28/2006

(preparation of phenylalanyl peptides as δ -opioid receptor ligands)

851883-43-7 HCAPLUS RN

L-Phenylalaninamide, 4-(aminocarbonyl)-L-phenylalanyl-D-alanylglycyl-N-(2-CN hydroxyethyl) $-N\alpha$ -methyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-A Ph Ме NH₂ H₂N Мe

PAGE 1-B

∕-он

L12 ANSWER 5 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

2005:295165 HCAPLUS ACCESSION NUMBER:

DOCUMENT NUMBER: 143:387341

TITLE: Short synthesis of C-terminal modified peptides by a

series-connection procedure

AUTHOR(S): Tian, Gui Jie; Qiu, Chuan Liang; Liu, Zhe; Wang, De

CORPORATE SOURCE: Institute of Materia Medica Chinese Academy of Medical

Sciences and Peking Union Medical College, Beijing,

100050, Peop. Rep. China

SOURCE: Chinese Chemical Letters (2005), 16(1), 31-34

CODEN: CCLEE7; ISSN: 1001-8417

PUBLISHER: Chinese Chemical Society

Journal DOCUMENT TYPE: LANGUAGE: English

CASREACT 143:387341 OTHER SOURCE(S): ΙT

866612-40-0 866612-41-1 866612-42-2

RL: RCT (Reactant); RACT (Reactant or reagent) (synthesis of C-terminal modified peptides)

866612-40-0 HCAPLUS RN

Glycinamide, L-tyrosylglycyl-L-phenylalanyl-N-(2-hydroxyethyl)- (9CI) (CA CN INDEX NAME)

RN 866612-41-1 HCAPLUS

CN Glycinamide, L-tyrosylglycyl-4-nitro-L-phenylalanyl-N-(2-hydroxyethyl)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 866612-42-2 HCAPLUS

CN β -Alaninamide, L-tyrosylglycyl-L-phenylalanyl-N-(2-hydroxyethyl)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT:

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 6 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

8

ACCESSION NUMBER: 2004:574827 HCAPLUS

DOCUMENT NUMBER: 141:271209

TITLE: Threonine at position 6 is not essential for the

immunosuppressive activity of HLA-DQ(β164-172)-

THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS

hexapeptide

AUTHOR(S): Stefanowicz, Piotr; Boratynski, Przemyslaw J.;

Staszewska, Anna; Wilczynski, Andrzej; Zimecki,

Michal; Szewczuk, Zbigniew

CORPORATE SOURCE: Faculty of Chemistry, University of Wroclaw, Wroclaw,

50-383, Pol.

SOURCE: Molecular Immunology (2004), 41(9), 911-917

CODEN: MOIMD5; ISSN: 0161-5890

PUBLISHER: Elsevier

10/551,771 08/28/2006

DOCUMENT TYPE: Journal LANGUAGE: English

IT 757967-17-2P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES

(Uses)

(immunosuppressive activity of HLA-DQ peptide analogs)

RN 757967-17-2 HCAPLUS

CN L-Tyrosinamide, L-arginylglycyl-L-α-aspartyl-L-valyl-N-(2-

hydroxyethyl) - (9CI) (CA INDEX NAME)

Absolute stereochemistry.

$$H_2N$$
 H_1
 H_2N
 H_3
 H_4
 H_5
 H_5
 H_5
 H_7
 H_7
 H_8
 H_8
 H_8
 H_8
 H_8
 H_9
 $H_$

REFERENCE COUNT: 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 7 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:549673 HCAPLUS

DOCUMENT NUMBER: 141:106735

TITLE: A solid phase method for synthesis of

peptide-spacer-lipid conjugates and preparation of

target liposome containing the conjugates

INVENTOR(S): Wu, Shih-Kuan; Chang, Ting-Kuang; Tseng, Chin-Lu;

Chen, Li-Rong; Shih, Kai-Hsiang

PATENT ASSIGNEE(S): Biotech Development Center, Taiwan SOURCE: Jpn. Kokai Tokkyo Koho, 82 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE

JP 2004189617 A2 20040708 JP 2002-355885 20021206
PRIORITY APPLN. INFO.: JP 2002-355885 20021206

OTHER SOURCE(S): CASREACT 141:106735

IT 632357-22-3DP, resin-bound 632357-23-4DP, resin-bound

632357-25-6DP, resin-bound

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(solid phase method for synthesis of peptide-spacer-lipid conjugates

and preparation of target liposome containing the conjugates)

RN 632357-22-3 HCAPLUS

CN L-Lysinamide, glycyl-1-(triphenylmethyl)-L-histidyl-N6-[(1,1-dimethylethoxy)carbonyl]-N-(2-hydroxyethyl)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 632357-23-4 HCAPLUS

CN L-Cysteinamide, S-[(acetylamino)methyl]-L-cysteinyl-L-methionyl-1- (triphenylmethyl)-L-histidyl-L-isoleucylglycyl-O-(1,1-dimethylethyl)-L-seryl-L-leucyl-L- α -aspartyl-O-(1,1-dimethylethyl)-L-seryl-O-(1,1-dimethylethyl)-L-threonyl-S- [(acetylamino)methyl]-N-(2-hydroxyethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-A

PAGE 1-B

RN 632357-25-6 HCAPLUS

CN L-Serinamide, glycyl-N5-[[[(2,3-dihydro-2,2,4,6,7-pentamethyl-5-benzofuranyl)sulfonyl]amino]iminomethyl]-L-ornithylglycyl-L- α -aspartyl-O-(1,1-dimethylethyl)-N-(2-hydroxyethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

L12 ANSWER 8 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:17423 HCAPLUS

DOCUMENT NUMBER: 140:72925

TITLE: Characterization and drug screening use of

phosphoinositolglycan-binding protein from plasma

10/551,771 08/28/2006

membrane of adipocytes

INVENTOR(S): Mueller, Guenter; Frick, Wendelin; Schneider, Rudolf;

Petry, Stefan; Urmann, Matthias

PATENT ASSIGNEE(S): Aventis Pharma Deutschland G.m.b.H., Germany

SOURCE: Eur. Pat. Appl., 41 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PATENT NO.					KIN	KIND DATE			APPLICATION NO.						DATE			
	EP	1378						2004	0107		EP	2002-	1504				0020	705	
		R:										IT,					MC,	PT,	
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		2490				AA						2003-							
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												EE,							
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			ΙE,	SI,	LT,	LV,	FI,	RO,	MK,	CY,	ΑI	, TR,	BG,	CZ,	EE,	ΗU,	SK		
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	CN	2006	903			Α		2006	0816		CN	2004- 2006-	1005	7471		2	0030	626	
	US	2004	2292	78		A1		2004	1118		US	2003-	4706	06		2	0030	703	
		7049				В2		2006	0523										
		2005						2005				2005-					0050	204	
	US	2006	1601	42		A1		2006	0720		US	2006-	3775	31		2	0060	316	
PRIO		Y APP									ΕP	2002-	1504	7		A 2	0020	705	
												2003-							
												2003-					0030		
											US	2003-	4706	06		A3 2	0030	703	
TT	611	1270-	20 7	n													-		

IT 640279-30-7P

RL: BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(PBP ligand; characterization and drug screening use of phosphoinositolglycan-binding protein (PBP) from plasma membrane of adipocytes)

RN 640279-30-7 HCAPLUS

CN L-Aspartamide, L-tyrosyl-L-methionyl-N1-(2-hydroxyethyl)- (9CI) (CA INDEX NAME)

10/551,771 08/28/2006

REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 9 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:971712 HCAPLUS

DOCUMENT NUMBER: 140:31479

TITLE: Solid phase method for preparation of peptide-lipid

conjugates for targeted liposome formulations INVENTOR(S): Wu, Shih-Kwang; Chang, Ting-Gung; Tseng, Chin-Lu;

Chen, Li-Jung; Shih, Kae-Shyang

PATENT ASSIGNEE(S): Development Center for Biotechnology, Taiwan

SOURCE: U.S. Pat. Appl. Publ., 19 pp., Cont.-in-part of U.S.

Pat. Appl. 2003 229,013.

CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2003229017	A1	20031211	US 2002-308644	20021203
US 2003229013	A1	20031211	US 2001-16569	20021203
CA 2413629	AA	20030607	CA 2002-2413629	20021205
CN 1453293	Α	20031105	CN 2002-155769	20021209
PRIORITY APPLN. INFO.:			US 2001-16569	A2 20011207

IT 632357-22-3DP, polymer-bound 632357-23-4DP, polymer-bound 632357-25-6DP, polymer-bound

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(solid phase method for preparation of peptide-lipid conjugates for targeted liposome formulations)

RN 632357-22-3 HCAPLUS

CN L-Lysinamide, glycyl-1-(triphenylmethyl)-L-histidyl-N6-[(1,1-dimethylethoxy)carbonyl]-N-(2-hydroxyethyl)- (9CI) (CA INDEX NAME)

RN 632357-23-4 HCAPLUS

CN L-Cysteinamide, S-[(acetylamino)methyl]-L-cysteinyl-L-methionyl-1- (triphenylmethyl)-L-histidyl-L-isoleucylglycyl-O-(1,1-dimethylethyl)-L-seryl-L-leucyl-L- α -aspartyl-O-(1,1-dimethylethyl)-L-seryl-O-(1,1-dimethylethyl)-L-threonyl-S- [(acetylamino)methyl]-N-(2-hydroxyethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

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RN 632357-25-6 HCAPLUS

CN L-Serinamide, glycyl-N5-[[[(2,3-dihydro-2,2,4,6,7-pentamethyl-5-benzofuranyl)sulfonyl]amino]iminomethyl]-L-ornithylglycyl-L- α -aspartyl-O-(1,1-dimethylethyl)-N-(2-hydroxyethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

L12 ANSWER 10 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:509780 HCAPLUS

DOCUMENT NUMBER:

139:392495

TITLE:

New proctolin analogues modified in position 2 and 5 of the peptide chain and their biological evaluation

10/551,771 08/28/2006

in insects

AUTHOR(S): Szeszel-Fedorowicz, Wioletta; Rosinski, Grzegorz;

Issberner, Jonathan; Osborne, Richard; Konopinska,

Danuta

CORPORATE SOURCE: Faculty of Chemistry University of Wroclaw, Wroclaw,

50-383, Pol.

SOURCE: Peptides 2000, Proceedings of the European Peptide

Symposium, 26th, Montpellier, France, Sept. 10-15, 2000 (2001), Meeting Date 2000, 849-850. Editor(s): Martinez, Jean; Fehrentz, Jean-Alain. Editions EDK:

Paris, Fr.

CODEN: 69EDWK; ISBN: 2-84254-048-4

DOCUMENT TYPE: Conference LANGUAGE: English

IT 395641-36-8

RL: AGR (Agricultural use); BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)

(proctolin analogs and their biol. activity in insects) -

RN 395641-36-8 HCAPLUS

CN L-Prolinamide, L-arginyl-L-tyrosyl-L-leucyl-N-(3-hydroxypropyl)- (9CI)

(CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

$$H_2N$$
 H_2N
 H_2N
 H_3N
 H_4N
 H_5
 H_5
 H_5
 H_6
 H_7
 H_7

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 11 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2001:386982 HCAPLUS

DOCUMENT NUMBER: 136:148140

TITLE: Myotropic effects of new proctolin analogues modified

in the position 5 of peptide chain in insects

AUTHOR(S): Szeszel-Fedorowicz, Wioletta; Rosinski, Grzegorz; Issberner, Jonathan; Osborne, Richard; Sliwowska,

Joanna; Konopinska, Danuta

CORPORATE SOURCE: Faculty of Chemistry, University of Wroclaw, Wroclaw,

PL 50-383, Pol.

SOURCE: Polish Journal of Pharmacology (2001), 53(1), 31-38

CODEN: PJPAE3; ISSN: 1230-6002

PUBLISHER: Polish Academy of Sciences, Institute of Pharmacology

DOCUMENT TYPE: Journal LANGUAGE: English

10/551,771 08/28/2006

395641-36-8P ΙT RL: BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation) (myotropic effects of proctolin analogs in insects) 395641-36-8 HCAPLUS RN L-Prolinamide, L-arginyl-L-tyrosyl-L-leucyl-N-(3-hydroxypropyl)- (9CI) CN (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

$$H_2N$$
 H_2
 H_2N
 H_3
 H_4
 H_5
 H_5
 H_5
 H_6
 H_7
 H_8
 $H_$

REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 12 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1965:403541 HCAPLUS

DOCUMENT NUMBER: 63:3541 ORIGINAL REFERENCE NO.: 63:671g-h

Gramicidin A. VI. The synthesis of valine- and TITLE:

isoleucine-gramicidin A

AUTHOR(S): Sarges, Reinhard; Witkop, Bernhard

CORPORATE SOURCE: U.S. Dept. of Health, Educ., & Welfare, Bethesda, MD Journal of the American Chemical Society (1965) 2020-7 SOURCE:

CODEN: JACSAT; ISSN: 0002-7863

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 63:3541

884483-21-0, Tryptophanamide, L-valyl-D-valyl-L-tryptophyl-Dleucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-884483-22-1, Tryptophanamide, D-valyl-L-tryptophyl-D-leucyl-Ltryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-885119-87-9, Tryptophanamide, D-valyl-L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-D-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-885119-89-1, Tryptophanamide, L-tryptophyl-D-leucyl-L-tryptophyl-Dleucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-

(preparation of)

884483-21-0 HCAPLUS CN Tryptophanamide, L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-Dleucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

Absolute stereochemistry.

RN

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RN 884483-22-1 HCAPLUS

CN Tryptophanamide, D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

PAGE 1-B

RN 885119-87-9 HCAPLUS

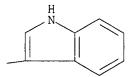
CN Tryptophanamide, D-valyl-L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-D-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

PAGE 1-B

RN 885119-89-1 HCAPLUS

CN Tryptophanamide, L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

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L12 ANSWER 13 OF 13 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1965:403540 HCAPLUS

DOCUMENT NUMBER: 63:3540
ORIGINAL REFERENCE NO.: 63:671d-g

TITLE: Gramicidin A. V. The structure of valine- and

isoleucine-gramicidin A

AUTHOR(S): Sarges, Reinhard; Witkop, Bernhard

CORPORATE SOURCE: U.S. Dept. of Health, Educ., & Welfare, Bethesda, MD

SOURCE: Journal of the American Chemical Society (1965),

87(9), 2011-20

CODEN: JACSAT; ISSN: 0002-7863

DOCUMENT TYPE: Journal LANGUAGE: English

IT 884483-21-0, Tryptophanamide, L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-

884483-22-1, Tryptophanamide, D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-885119-87-9, Tryptophanamide, D-valyl-L-valyl-D-valyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-885119-89-1, Tryptophanamide, L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L-(preparation of)

RN 884483-21-0 HCAPLUS

CN Tryptophanamide, L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-

leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

RN 884483-22-1 HCAPLUS

CN Tryptophanamide, D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

RN 885119-87-9 HCAPLUS

CN Tryptophanamide, D-valyl-L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-D-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

PAGE 1-B

PN. 865119-89-1 HCAPLUS

Pryptophanamide, L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

RN 885119-87-9 HCAPLUS

CN Tryptophanamide, D-valyl-L-valyl-D-valyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-D-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

PAGE 1-B

RN 885119-89-1 HCAPLUS

CN Tryptophanamide, L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-L-tryptophyl-D-leucyl-N-2-hydroxyethyl-, L- (7CI) (CA INDEX NAME)

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